

| COMPOUND IDENTIFIED   | 1.0 | 6.0 | 24.0 | 48.0 | 72.0 | 96.0 |
|---|-----|-----|------|------|------|------|
| Nonane, 4-methyl  |     | 1.9 |      |      |      |      |
| Octane, 3,6-dimethyl  |     | 2.3 | .    |      |      |      |
| Pentane, 2,2,4-trimethyl (Isooctane)                        | 1.6 |     |      | 0.7  |      |      |
| Pinene, a<br>(2,6,6-Trimethyl-bicyclo[3.1.1]hept-2-ene<br>) | 3.4 |     |      |      |      |      |
| Styrene   | 7.0 | 3.1 | 5.5  | 5.5  | 4.2  | 6.5  |
| Toluene (Methylbenzene)                                     | 9.2 | 0.7 |      |      |      |      |
| Unidentified  |     |     |      |      |      | 1.5  |
| Xylene (para and/or meta)                                   | 4.8 |     |      |      |      |      |
| Xylene, ortho   | 2.0 |     |      |      |      |      |
| cis-1-Ethyl-3-methykylohexane*                              | 4.0 |     |      |      |      |      |

\*Indicates best NIST/EPA/NIH library match only.

Individual compounds and TVOC (total volatile organic compounds) are calibrated relative to toluene.

TVOC detection limit: 0.9  $\mu\text{g}/\text{m}^3$ . Individual detection limits may vary, depending on instrument response.

### Cushion C

Environmental Chamber, SA3  
Product Loading: 0.42 m<sup>2</sup>/m<sup>3</sup>  
Test Conditions: 1.00 ACH  
50.0% RH ± 5.0% RH  
23.0 degs C ± 2.0 degs C  
Test Period: 06/22/95 - 06/26/95

#### CHAMBER TVOC CONCENTRATIONS FROM 0.000 TO 96.000 HOURS

| ELAPSED EXPOSURE HOUR | Tvoc CONCENTRATION ug/m <sup>3</sup> |
|-----------------------|--------------------------------------|
| 0.000                 | 0.0                                  |
| 1 .000                | 915.0                                |
| 6.000                 | 826.6                                |
| 24.000                | 265.3                                |
| 48.000                | 366.9                                |
| 72.000                | 253.1                                |
| 96.000                | 38.8                                 |

#### IVOC CONCENTRATIONS (µg/m<sup>3</sup>) FROM 0.000 TO 96.000 HOURS

| COMPOUND IDENTIFIED                    | 1.0  | 6.0   | 24.0 | 48.0 | 72.0 | 96.0 |
|--|------|-------|------|------|------|------|
| 1,3-Dioxolan-2-one, cl-methyl'         |      | 5.9   | 4.8  | 11.0 | 9.1  |      |
| 1-Octanol, 3,7-dimethyl*               | 2.0  | 2.4   |      |      |      |      |
| I-Pentanol, 2-methyl-*                 | 1.5  |       |      |      |      |      |
| 1-Pentanol, 4-methyl-2-propyl- (9Cl)*  | 3.0  |       |      |      |      |      |
| 2,2,7,7-Tetramethyloctane*             | 15.6 | 5.5   |      |      |      |      |
| 2,6-Di-tert-butyl-4-methylphenol (BHT) | 66.9 | 106.1 | 31.3 | 82.5 | 72.1 | 38.8 |
| 2-Hexanol, 5-methyl-*                  | 4.5  |       |      |      |      |      |
| 2-Propanol (Isopropanol)               | 2.2  |       |      |      |      |      |

| COMPOUND IDENTIFIED  | 1.0   | 6.0   | 24.0 | 48.0 | 7i.o | 96.0 |
|--|-------|-------|------|------|------|------|
| <u>2-Propanol, 1 ,3-dichloro-</u>  | 74.8  | 106.3 | 41.6 | 65.2 | 86.3 |      |
| <u>2-Propanone, 1-(dimethylamino)-*</u>  | 5.6   |       |      |      |      |      |
| <u>2H-2,4a-Methanonaphthalene,<br/>1,3,4,5,6,7-hexahydro-1,1,5,5-tetramethyl<br/>-, (2S)-*</u> | 1.6   | 1.3   | 1.0  | 1.7  | 0.8  |      |
| <u>4-Octanol, 4,7-dimethyl-*</u>   |       | 1.2   |      |      |      |      |
| <u>Acetamide, N,N-dimethyl- (8Cl9Cl)</u>   |       | 82.0  | 65.0 | 61.8 |      |      |
| <u>Acetone (2-Propanone)</u>   | 11.2  |       |      |      |      |      |
| <u>Benzene, 1-ethyl-4-methyl<br/>(4-Ethyltoluene)</u>  | 2.6   |       |      |      |      |      |
| <u>Benzene, 1-methyl-4-(1-methylethyl)<br/>(p-Cymene; 4-Isopropyltoluene)*</u>                 | 2.2   |       |      |      |      |      |
| <u>Benzene,(1 ,1-Dimethylethyl)- .<br/>(t-Butylbenzene)*</u>                                   | 6.8   | 2.0   |      |      |      |      |
| <u>Butane,<br/>2,2'-[methylenebis(oxy)]bis[2-methyl-*</u>                                      | 1.5   |       |      |      |      |      |
| <u>Cyclohexane, propyl</u>   | 1.7   |       |      |      |      |      |
| <u>Cyclohexanone, 4-methyl-*</u>   | 1.4   |       |      |      |      |      |
| <u>Cyclohexasiloxane, dodecamethyl*</u>  | 8.9   | 9.4   | 3.1  | 1.6  |      |      |
| <u>Cyclopentasiloxane, decamethyl*</u>   | 145.4 | 123.7 | 7.3  | 0.8  |      |      |
| <u>Cyclotetrasiloxane, octamethyl</u>  | 132.3 | 32.1  |      |      |      |      |
| <u>Decane, 2,5,9-trimethyl- (9Cl)</u>  | 5.0   | 4.6   | 0.7  |      |      |      |
| <u>Decane, 2-methyl*</u>   | 5.1   | 3.4   |      |      |      |      |
| <u>Decane, 3-methyl</u>  | 16.8  | 5.1   |      |      |      |      |
| <u>Decane, 4-methyl</u>  | 10.3  | 6.1   |      |      |      |      |
| <u>Dodecane</u>  | 3.4   | 1.2   |      |      |      |      |
| <u>Ethane, 1,1, 1-trkhloro</u>   | 2.8   |       |      |      |      |      |
| <u>Ethane, 1,2-dichloro</u>  | 1.9   |       |      |      |      |      |
| <u>Ethane,1,1'oxybis[2-chloro-<br/>(s-Dichloroethyl ether)</u>                                 |       | 1.2   |      |      |      |      |
| <u>Ethanol</u>   | 10.9  |       |      |      |      |      |

| COMPOUND IDENTIFIED                                  | 1.0  | 6.0   | 24.0 | 48.0  | 72.0 | 96.0 |
|--|------|-------|------|-------|------|------|
| Ethene, 1,1,2,2-tetrachloro<br>(Tetrachloroethylene) | 2.7  |       |      |       |      |      |
| Formamide, N,N-dimethyl                              | 2.8  | 78.9  | 13.5 | 6.1   |      |      |
| Hepiane, 2,2,4,6,6-pentamethyl (8Cl9Cl)*             | 6.8  | 1.9   |      |       |      |      |
| Heptane, 2,2,6,6-tetramethyl*                        | 11.2 | 1.3   |      |       |      |      |
| Heptane, 4-ethyl-2,2,6,6-tetramethyl-<br>(9Cl)       | 18.9 | 7.3   |      |       |      |      |
| Hexanal (Hexaldehyde)                                | 1.4  |       |      |       |      |      |
| Hexane, 1 ,1'-Oxybis (Hexyl ether)*                  | 1.5  |       |      |       |      |      |
| Hexane, 2,2,4-trimethyl                              | 2.3  |       |      |       |      |      |
| Isooctane, (ethenyoxy)- (9Cl)*                       |      | 1.3   |      |       |      |      |
| Methanol   | 28.7 |       | 5.4  |       |      |      |
| Methylene chloride (Dichloromethane)                 | 5.1  |       |      |       |      |      |
| N,N-Dimethyl-2-propenamide                           | 5.0  | 100.9 | 70.9 | 126.8 | 82.6 |      |
| Nonane   | 2.9  |       |      |       |      |      |
| Nonane, 2,5-dimethyl-*                               | 13.1 | 5.9   |      |       |      |      |
| Nonane, 3,7-dimethyl*                                | 3.4  |       |      |       |      |      |
| Nonane, 3-ethyl                                      | 1.9  |       |      |       |      |      |
| Nonane, 3-methyl                                     | 1.5  |       |      |       |      |      |
| Nonane, 4,5-dimethyl*                                | 2.1  | -     |      |       |      |      |
| Octane, 2,2-dimethyl*                                | 4.7  |       |      |       |      |      |
| Octane, 2,3,6-trimethyl-*                            | 2.7  |       |      |       |      |      |
| Octane, 2,3-dimethyl*                                | 2.3  |       |      |       |      |      |
| Octane, 2,6-dimethyl                                 | 3.0  |       |      |       |      |      |
| Oxetane, 2,4-dimethyl-, trans-*                      | 5.9  | 1.2   |      |       |      |      |
| Oxirane, (chloromethyl)-•                            |      |       |      | 1.4   |      |      |
| Pentane*   | 1.3  |       |      |       |      |      |
| Pentane, 2,2,4-trimethyl (Isooctane)                 | 9.2  | 1.4   |      |       |      |      |
| Pentane, 2,3,4-trimethyl                             | 3.4  |       |      |       |      |      |

| COMPOUND IDENTIFIED   | 1.0   | 6.0   | 24.0 | 48.0 | 72.0 | 96.0 |
|---|-------|-------|------|------|------|------|
| Pentane, 3-ethyl-2,2-dimethyl- (8CI9Cl)*                        | 2.8   |       |      |      |      |      |
| Phenol  |       | 2.7   |      | 1.5  | 2.0  |      |
| Propane,? ,2,3-trichloro  | 129.0 | 111.5 | 20.6 | 6.5  |      |      |
| Toluene (Methylbenzene)   | 59.9  | 2.5   |      |      |      |      |
| Undecane  | 9.4   | 4.1   |      |      |      |      |
| Undecane, 6-methyl-*  | 3.2   | 1.3   |      |      |      |      |
| Unidentified  | 5.7   | 2.3   |      |      |      |      |
| Xylene (para and/or meta)                                       | 7.1   |       |      |      |      |      |
| Xylene, ortho   | 2.8   |       |      |      |      |      |
| a-Methylstyrene (iso-Propenylbenzene; (I-Methylethenyl)benzene) | 3.7   | 2.3   |      |      |      |      |

\*indicates best NIST/EPA/NIH library match only.

Individual compounds and TVOC (total volatile organic compounds) are calibrated relative to toluene.

TVOC detection limit: 0.9 µg/m<sup>3</sup>. Individual detection limits may vary, depending on instrument response.

## Cushion D

Environmental Chamber. SA4  
Product Loading: 0.40 m<sup>2</sup>/m<sup>3</sup>  
**Test** Conditions: 1.00 ACH  
50.0% RH ± 5.0% RH  
23.0 degs C ± 2.0 degs C  
**Test Period:** 06/26/95 - 06/30/95

### CHAMBER TVOC CONCENTRATIONS FROM 0.000 TO 96.000 HOURS

| ELAPSED EXPOSURE HOUR | TVOC CONCENTRATION (µg/m <sup>3</sup> ) |
|-----------------------|---|
| 0.000                 | 0.0                                     |
| 1 .000                | 349.0                                   |
| 6.000                 | 634.9                                   |
| 24.000                | 421.6                                   |
| 48.000                | 345.8                                   |
| 72.000                | NA*                                     |
| 96.000                | 212.0                                   |

\*NA: data not available due to laboratory error.

### IVOC CONCENTRATIONS (µg/m<sup>3</sup>) FROM 0.000 TO 96.000 HOURS

| COMPOUND IDENTIFIED  | 1.0  | 6.0   | 24.0  | 48.0  | 96.0  |
|--|------|-------|-------|-------|-------|
| 1,4-Dimethylpiperazine   | 58.5 | 319.1 | 71.4  | 55.3  |       |
| 1H-Cyclopropa[a]naphthalene,<br>1a,2,3,5,6,7,7a,7b-octahydro-1,1,7,7a<br>-tetramethyl-,[1aR-(1aα,7α,7aα,7bα)]<br>* |      | 1.3   | 1.7   |       | 1.3   |
| 2,2,7,7-Tetramethyloctane*   | 0.8  |       |       |       |       |
| 2,6-Di-tert-butyl-4-methylphenol<br>(BHT)  | 23.5 | 108.1 | 149.9 | 122.1 | 100.9 |
| 2-Butenal*   | 2.1  |       |       |       |       |

| COMPOUND IDENTIFIED                      | 1.0  | 6.0  | 24.0 | 48.0 | 96.0 |
|--|------|------|------|------|------|
| 2-Propanol (isopropanol)                 | 2.5  |      |      |      |      |
| 4-Isobutylmorpholine*                    | 1.1  |      |      |      |      |
| Acetamide, N,N-dimethyl-(8Cl9Cl)         | 13.0 | 19.4 | 15.7 | 11.0 | 7.4  |
| Acetate, butyl                           | 1.1  |      |      |      |      |
| Acetic acid, hydroxy-•                   |      |      |      | 1.0  |      |
| Acetone (2-Propanone)                    | 5.9  |      |      |      |      |
| Benzene, 1,2,4-trimethyl                 | 0.7  |      |      |      |      |
| Benzene, 1,2-dichloro                    | 7.7  | 8.0  | 5.3  | 3.8  | 1.9  |
| Benzene, ethyl                           | 1.0  |      |      |      |      |
| Benzene, propyl                          | 1.0  |      |      |      |      |
| Butane, 2-methyl (Isopentane)            | 1.8  |      |      |      |      |
| Cyclohexasiloxane, dodecamethyl*         | 0.8  | 1.7  | 2.3  |      |      |
| Cyclopentasiloxane, decamethyl*          | 3.8  | 2.8  |      |      |      |
| Decane                                   | 3.6  |      |      |      |      |
| Decane, 2-methyl                         | 2.5  |      |      |      |      |
| Decane, 5-methyl*                        | 0.9  |      |      |      |      |
| Dodecane                                 |      | 1.3  |      |      |      |
| Ethane, 1,1,1-trichloro                  | 38.5 | 23.6 | 14.0 | 10.0 | 5.6  |
| Ethanol                                  | 2.2  |      |      |      |      |
| Ethanol, 2-butoxy                        | 1.8  | 2.0  |      |      |      |
| Heptane, 2,2,4,6,6-pentamethyl (8Cl9Cl)* | 2.0  |      |      |      |      |
| Hexane, 2,2,4-trimethyl*                 | 0.5  |      |      |      |      |
| Nonane                                   | 0.8  |      |      |      |      |
| Nonane, 3,7-dimethyl*                    | 1.5  |      |      |      |      |
| Nonane, 4-methyl                         | 1.2  |      |      |      |      |
| Octane, 2,2,6-trimethyl*                 | 0.9  |      |      |      |      |
| Octane, 2,3,6-trimethyl-*                | 0.8  |      |      |      |      |

| COMPOUND IDENTIFIED                                     | 1.0   | 6.0   | 24.0  | 48.0  | 96.0 |
|---|-------|-------|-------|-------|------|
| Octane, 2,3-dimethyl                                    | 1.2   |       |       |       |      |
| Pentane, 2,2,4-trimethyl (Isooctane)                    | 2.1   |       |       |       |      |
| Pentane, 2,3,4-trimethyl*                               | 0.6   |       |       |       |      |
| Phenol  |       | 1.0   |       |       |      |
| Pinene, a<br>(2,6,6-Trimethyl-bicyclo[3.1.1]hept-2-ene) | 1.1   |       |       |       |      |
| Styrene   | 40.6  | 40.3  | 23.1  | 12.2  | 4.2  |
| Toluene (Methylbenzene)                                 | 16.9  | 4.0   |       |       |      |
| Undecane  | 1.6   | 1.0   |       |       |      |
| Unidentified  | 100.8 | 101.2 | 138.2 | 130.4 | 90.7 |
| Xyiene, ortho   | 1.5   |       |       |       |      |

\*Indicates best NIST/EPA/NIH library match only.

Individual compounds and TVOC (total volatile organic compounds) are calibrated relative to toluene.

TVOC detection limit: 0.9 µg/m<sup>3</sup>. Individual detection limits may vary, depending on instrument response.

### Cushion E

Environmental Chamber: SA5  
Product Loading: 0.42 m<sup>2</sup>/m<sup>3</sup>  
Test Conditions: 1.00 ACH  
50.0% RH ± 5.0% RH  
23.0 degs C ± 2.0 degs C  
Test Period: 06/26/95 - 06/30/95

#### CHAMBER TVOC CONCENTRATIONS FROM 0.000 TO 96.000 HOURS

| ELAPST VOC<br>EXPOSURE/CONCENTRATION<br>HOU(µg/m <sup>3</sup> ) |        |
|---|--------|
| 0.000   | 0.0    |
| 1 .000  | 1447.1 |
| 6.000   | 1631.1 |
| 24.000  | 1260.0 |
| 48.000  | 856.1  |
| 72.900  | 572.9  |
| 96.000  | 177.7  |

#### IVOC CONCENTRATIONS (µg/m<sup>3</sup>) FROM 0.000 TO 96.000 HOURS

| COMPOUND IDENTIFIED                                       | 1.0 | 6.0  | 24.0 | 48.0 | 72.0 | 96.0 |
|---|-----|------|------|------|------|------|
| 1,2-Propanediol (Propylene glycol)                        | 3.1 | 27.1 | 28.1 | 26.8 | 20.2 |      |
| 1,3-Butanediol*   | 4.2 | 44.9 | 40.4 | 41.5 | 30.6 |      |
| 1,3-Cyclohexadiene,<br>1-methyl-4-(1-methylethyl)-*       | 2.5 |      | 2.2  |      |      |      |
| 1 ,4-Cyclohexadiene,<br>1-methyl-4-(1-methylethyl)- (9CI) |     |      | 0.8  |      |      |      |
| 1,4-Dimethylpiperazine                                    |     |      |      | 2.1  |      |      |
| 1,4-Dioxane,<br>2-methyl-3-methylene-*                    |     |      |      | 3.1  |      |      |
| 1-Heptanol, 6-methyl*                                     | 1.7 |      |      |      |      |      |

| COMPOUND IDENTIFIED  | '1.0   | 6.0    | 24.0   | 48.0  | 72.0  | 96.0  |
|--|--------|--------|--------|-------|-------|-------|
| 1-Undecene, 8-methyl-*   | 6.5    | 5.3    |        |       |       |       |
| 2(3H)-Furanone, dihydro<br>(Butyrolactone)*                          |        | 3.3    | 3.5'   | 2.9   | 1.9   |       |
| 2,3-Butanediol   | 2.9    | 28.2   | 24.5   | 26.2  | 19.8  |       |
| 2,6-Octadien-1-ol, 3,7-dimethyl-,<br>(Z)-*                           |        | 1.7    |        |       |       |       |
| 2-Butanone, 3-hydroxy-*  |        | 2.1    |        |       |       |       |
| 2-Decene, 4-methyl-, (Z)*  | 6.2    |        |        |       |       |       |
| 2-Decene, 7-methyl-, (Z)-*   | 5.7    | 2.7    |        |       |       |       |
| 2-Decene, 8-methyl-, (Z)-*   | 3.0    | 1.6    |        |       |       |       |
| 3-Cyclohexene-1-methanol,<br>$\alpha,\alpha,4$ -trimethyl-, acetate* | 43.7   | 69.1   | 49.8   | 54.7  | 35.7  | 20.2  |
| 4-Dodecene*  | 8.0    | 6.9    |        |       |       |       |
| 4-Phenylcyclohexene  | 2.3    | 9.8    | 8.4    | 5.5   | 3.0   | 1.6   |
| 4-Undecene, (Z)-*  | 7.0    | 6.3    |        |       |       |       |
| 5-Undecene, (E)  | 8.3    | 6.4    |        |       |       |       |
| 6-Dodecene, (E)- (8Cl9Cl)*   |        | 3.1    |        |       |       |       |
| Acetamide, N,N-dimethyl-<br>(8Cl9Cl)                                 |        | 2.7    | 2.5    | 3.4   | 1.7   |       |
| Acetate, methyl (Acetic acid,<br>Methyl ester)                       | 11.4   | 1.7    |        |       |       |       |
| Acetic acid  | 1170.4 | 1271.3 | 1004.8 | 641.7 | 426.1 | 137.5 |
| Acetic acid, phenylmethyl ester<br>(Benzyl acetate)                  |        | 8.4    | 3.9    | 2.1   | 1.2   |       |
| Acetone (2-Propanone)  | 16.8   | 3.7    |        |       |       |       |
| Benzaldehyde   |        | 8.1    | 4.1    | 2.8   | 1.7   | 2.2   |
| Benzyl alcohol<br>(Benzenemethanol)                                  |        | 7.1    | 4.4    | 6.0   | 3.8   | 1.6   |
| Bicyclo[3.1.1]heptane-2-methano<br>1,6,6-dimethyl-, acetate*         | 0.3    |        |        |       |       |       |

| COMPOUND IDENTIFIED   | 1.0  | 6.0  | 24.0 | 48.0 | 72.0 | 96.0 |
|---|------|------|------|------|------|------|
| Bicyclo[7.2.0]undec-4-ene,<br>4,11,11-trimethyl-8-methylene-, [1<br>R-(1R@,4Z,9S@)]-* | 1.2  | 1.4  | 1.3  |      |      |      |
| Cedrene, $\alpha^*$   | 5.8  | 5.3  | 5.7  | 2.5  | 4.8  | 1.7  |
| Cyclohexane, 1,1'-oxybis*   |      | 2.5  | 2.1  | 1.2  |      |      |
| Cyclohexanol  |      | 1.8  |      |      |      |      |
| Cyclohexanol,<br>1-methyl-4-(1-methylethenyl)-,<br>acetate*                           | 1.2  | 1.3  | 0.9  |      |      |      |
| Cyclohexene,<br>1-methyl-4-(1-methylethylidene)-*                                     | 22.3 | 20.1 | 9.0  | 2.7  | 1.0  | 0.9  |
| Cyclopentasiloxane, decamethyl*   | 21.4 | 2.4  |      |      |      |      |
| Cyclopropane, 1-ethyl-2-heptyl-<br>(9CI)*   | 7.3  | 4.7  |      |      |      |      |
| Cyclotetrasiloxane, octamethyl  | 1.3  |      |      |      |      |      |
| Decane  |      | 1.4  |      | 1.9  |      |      |
| Ethanol   | 3.1  | 1.8  |      |      |      |      |
| Limonene (Dipentene;<br>1-Methyl-4-(1-methylethyl)cyclohexene)                        | 25.9 | 35.0 | 31.0 | 13.9 | 6.0  | 6.6  |
| Methanol  | 24.6 |      |      |      |      |      |
| Nonane, 2,6-dimethyl*   | 2.0  |      |      |      |      |      |
| Propanoic acid  |      | 19.6 | 16.8 | 12.3 | 9.5  | 1.5  |
| Toluene (Methylbenzene)   | 4.8  |      |      |      |      |      |

| COMPOUND IDENTIFIED  | 1.0  | 6.0 | 24.0 | 48.0 | 72.0 | 96.0 |
|--|------|-----|------|------|------|------|
| <b>Tricyclo[5.4.0.0<sup>2,8</sup>]undec-9-ene,<br/>2,6,6,9-tetramethyl-*</b> | 7.7  | 8.3 | 7.7  | 4.8  | 4.8  | 3.1  |
| Unidentified   | 14.7 | 4.4 | 5.0  | 1.1  | 1.3  | 0.9  |

\*Indicates best NIST/EPA/NIH library match only.

Individual compounds and TVOC (total volatile organic compounds) are calibrated relative to toluene.

TVOC detection limit: 0.9  $\mu\text{g}/\text{m}^3$ . Individual detection limits may vary, depending on instrument response.

**Carpet B at Elevated Conditions (70 °C, 8% RH, high loading)**

**Environmental Chamber:** SC4  
**Product Loading:** 1.64 m<sup>2</sup>/m<sup>3</sup>  
**Test Conditions:** 1.00 ACH  
8.0% RH ± 2.0% RH  
70.0 degs C ± 5.0 degs C  
**Test Period:** 06/21/95 - 06/21/95

**CHAMBER TVOC CONCENTRATIONS FROM 0.000 TO 99.000 HOURS**

| ELAPSED EXPOSURE HOUR | Tvoc CONCENTRATION (µg/m <sup>3</sup> ) |
|-----------------------|---|
| 0.000                 | 9.5                                     |
| 1.000                 | > 35649.5                               |
| 6.000                 | 7127.5                                  |

**IVOC CONCENTRATIONS (µg/m<sup>3</sup>) FROM 0.000T06.000 HOURS**

| COMPOUND IDENTIFIED              | 0.0 | 1.0    | 6.0   |
|----------------------------------|-----|--------|-------|
| 1,2-Ethanediol (Ethylene glycol) |     | 98.3   | 466.0 |
| 1,3-Butanediol, (S)-*            |     |        | 4.3   |
| 1,4-Pentadien-3-ol*              |     | 36.3   | 7.9   |
| I-Butanol (N-Butylalcohol)       |     | 165.1  | 12.3  |
| 1-Decanol (N-Decyl alcohol)      |     | 60.6   | 21.6  |
| I-Decene, 8-methyl               |     | 292.6  | 92.5  |
| I-Dodecanol                      |     | 766.5  | 63.4  |
| 1-Dodecene*                      |     | 51.8   | 17.5  |
| 1-Heptanol, 6-methyl*            |     | 1163.7 | 288.2 |
| 1-Hexene                         |     | 4.5    |       |
| I-Octanol, 2-butyl- (8Cl9Cl)*    |     | 7.1    | 178.7 |
| 1-Octene, 2,6-dimethyl-*         |     | 74.6   | 22.5  |

| COMPOUND IDENTIFIED  | 0.0 | 1.0    | 6.0   |
|--|-----|--------|-------|
| I-Propanol ( <u>Propyl alcohol</u> )                                       |     | 57.7   | 13.3  |
| I-Propanol, 2-(2-hydroxypropoxy)*  |     | 405.1  | 371.5 |
| 1-Propene, 3-chloro-2-methyl-*   |     | 9.1    | 2.9   |
| 1-Tetradecanol*  |     | 468.6  | 13.8  |
| 2(1 <i>i</i> -i)-Naphthalenone,<br>octahydro-4a,7,7-trimethyl-,<br>trans-• |     | 142.4  | 9.1   |
| 2,2,4-Trimethyl-1,3-pentanediol<br>monoisobutyrate (Texanol)               |     | 267.2  | 33.7  |
| 2,2-Dimethyl-1-isopropyl-1,3-prop<br>anediol monoisobutyrate<br>(Texanol)  |     | 137.3  | 29.0  |
| 2,6-Di- <i>tert</i> -butyl-4-methylphenol<br>(BHT)                         |     | 2351.0 | 328.3 |
| 2,6-Octadien-1-ol, 3,7-dimethyl-,<br>acetate, (Z)-*                        |     | 78.8   | 6.3   |
| 2,6-Octadien-1-ol,3,7-dimethyl, (E)<br>- (Geraniol)*                       | •   | 10.4   | 6.4   |
| 2-Butanone (Methyl ethyl ketone,<br>MEK)                                   |     | 5.0    |       |
| 2-Butenal, 2-ethenyl-*   |     |        | 2.4   |
| 2-Cyclohexen-1-one,<br>2-methyl-5-(1-methylethenyl)-*                      |     | 14.3   | 6.9   |
| 2-Decene, g-methyl-, (Z)-*   |     | 217.8  | 44.4  |
| 2-Decene, (Z)-*  |     |        |       |
| 2-Decene, 3-methyl-, (Z)-*   |     | 327.7  | 79.0  |
| 2-Decene, 5-methyl-, (Z)-*   |     | 387.6  | 173.0 |
| 2-Decene, 7-methyl-, (Z)-*   |     | 220.0  | 35.8  |
| 2-Dodecen-1-yl(-)succinic<br>anhydride*                                    |     | 1090.4 |       |
| 2-Dodecene, (E)*   |     | 322.3  |       |
| 2-Naphthalenemethanol,<br>α-methyl-, (+/-)-*                               |     |        | 6.3   |

| COMPOUND IDENTIFIED  | 0.0 | 1.0    | 6.0   |
|--|-----|--------|-------|
| 2-Octene, (E)-*  |     | 12.6   |       |
| 2-Octene, (Z)-*  |     | 8.7    |       |
| 2-Octene, 4-ethyl-*  |     | 96.7   |       |
| 2-Pentanone, 4-hydroxy-4-methyl-<br>(8CI9Cl)*                        |     | 64.6   | 4.9   |
| 2-Pentanone, 4-methyl (Methyl<br>isobutyl ketone, MIBK)*             |     | 5.2    |       |
| 2-Propanol (Isopropanol)   |     | 21.1   | 4.1   |
| 2-Propen-1-ol, 2-methyl  |     | 6.0    | 0.8   |
| 2-Propenal*  |     |        | 3.9   |
| 2-Propenoic acid, methyl ester*                                      |     | 37.6   | 13.6  |
| 2-Undecene, 2,5-dimethyl-*   |     | 500.0  | 172.2 |
| 2H-Benzocyclohepten-2-one,<br>decahydro-4a-methyl-, trans-•          |     | 1214.1 |       |
| 3,5-Cyclohexadiene-1,2-dione,<br>3,5-bis(1, 1-dimethylfethyl)-•      |     |        | 11.5  |
| 3-Buten-2-ol*  |     | 3.9    |       |
| 3-Cyclohexene-1-methanol,<br>$\alpha,\alpha,4$ -trimethyl-*          |     | 82.8   | 29.5  |
| 3-Cyclohexene-1-methanol,<br>$\alpha,\alpha,4$ -trimethyl-, acetate* |     | 61.5   | 9.4   |
| 3-Decene*  |     | 273.3  | 106.1 |
| 3-Decene, 2-methyl-, (Z)-•   |     | 985.6  | 281.6 |
| 3-Dodecene, (E)*   |     | 527.3  | 158.6 |
| 3-Dodecene, (Z)- (8CI9Cl)*   |     | 63.6   |       |
| 3-Ethyl-2-methyl-1-heptene*  |     |        | 3.1   |
| 3-Octene, 2,6-dimethyl-*   |     | 99.1   | 18.2  |
| 3-Undecene, 2-methyl-, (E)-•   |     | 1660.4 | 469.6 |
| 4-Decene, 3-methyl-, (E)-•   |     | 697.0  |       |
| 4-Dodecene*  |     | 521.5  | 152.5 |

| COMPOUND IDENTIFIED  | 0.0 | 1.0                  | 6.0   |
|--|-----|----------------------|-------|
| <b>4-Isobutylmorpholine</b>  |     | <b>&gt; 1317.7**</b> | 547.3 |
| <b>4-Nonene, 3-methyl, (Z)</b>   |     | <b>551.7</b>         | 126.6 |
| <b>4-Nonene, 5-methyl-*</b>  |     | 56.0                 | 34.0  |
| <b>4-Octene, (E)*</b>  |     | 35.0                 |       |
| <b>4-Octene, 2,3,6-trimethyl-*</b>   |     |                      | 130.5 |
| 4-Phenylcyclohexene  |     | 58.7                 | 10.2  |
| <b>4-Piperidinecarboxylic acid, 4-phenyl-, ethyl ester*</b>                  |     | 633.1                | 26.2  |
| <b>4-tert-Butylcyclohexyl acetate*</b>                                       |     |                      | 3.3   |
| <b>5-Dodecene, (E)*</b>  |     | 191.6                | 49.5  |
| <b>5-Undecene, (E)*</b>  |     | 1045.9               | 331.1 |
| <b>5-Undecene, (Z)- (8Cl9Cl)*</b>  |     | 1789.0               | 403.0 |
| <b>6-Dodecene, (E)- (8Cl9Cl)</b>   |     | 816.4                | 348.0 |
| <b>Acetamide, N,N-dimethyl- (8Cl9Cl)</b>                                     |     | 404.3                | 60.5  |
| <b>Acetamide, N-methyl-N-[4-[4-methoxy-1-hexa hydropyridyl]-2-butynyl]-*</b> |     | 264.1                | 28.2  |
| Acetate, ethyl   |     | 4.0                  |       |
| Acetic acid  | 0.9 |                      | 153.3 |
| <b>Acetone (2-Propanone)</b>   |     | 78.0                 | 23.1  |
| <b>Azulene, 7-ethyl-1,4-dimethyl-*</b>                                       |     |                      | 8.1   |
| <b>Benzene, 1-ethyl-4-methyl (4-Ethyltoluene)</b>                            |     | 43.1                 |       |
| Benzene, ethyl   |     | 23.4                 |       |
| <b>Benzene.1 ,1'-Oxybis- (Diphenyl ether)'</b>                               |     | 126.4                | 10.3  |
| <b>Benzoic acid, 2-hydroxy-, 3-methylbutyl ester"</b>                        |     | 399.7                | 53.6  |
| <b>Benzoic acid, 2-hydroxy-, pentyl ester*</b>                               |     | 333.3                | 10.1  |

| COMPOUND IDENTIFIED   | 0.0 | 1.0   | 6.0   |
|---|-----|-------|-------|
| Benzoic acid,<br>3-(5-hydroxy-1-pentenyl)-, methyl ester, (E)-*         |     | 138.0 | 4.7   |
| Benzothiatole   |     |       | 8.8   |
| Butane  |     | 6.7   |       |
| Butanoic acid   |     |       | 10.3  |
| Caryophyllene oxide*  |     | 464.7 |       |
| Cyclohexane, (1-methylpropyl)-*   |     | 44.1  | 4.2   |
| Cyclohexane, 1,1,2-trimethyl*   |     | 11.6  | 1.2   |
| Cyclohexane,<br>1,1-dimethyl-2-propyl-*                                 |     | 278.0 | 163.0 |
| Cyclohexane,<br>1-ethenyl-3-methylene-5-(1-propenylidene)-*             |     |       | 2.0   |
| Cyclohexane, 1-ethyl-2-methyl-, cis- (8Cl9Cl)*                          |     | 17.5  | 3.9   |
| Cyclohexane, methyl   |     | 5.2   |       |
| Cyclohexane, octyl*   |     |       | 33.6  |
| Cyclohexane, propyl   |     | 61.0  |       |
| Cyclohexane, t-1-ethyl-4-methyl*  |     | 32.8  |       |
| Cyclohexanol,<br>2-methylene-3-(1-methylethyl)-, acetate, cis-•         |     | 52.6  | 15.1  |
| Cyclopentane, 1,2,4-trimethyl,<br>(1 $\alpha$ ,2 $\beta$ ,4 $\alpha$ )* |     | 11.0  | 1.7   |
| Cyclopentasiloxane, decamethyl*   |     | 99.5  | 25.9  |
| Decane  |     | 910.6 | 46.7  |
| Decane, 6-ethyl-2-methyl- (9Cl)*  |     | 12.1  |       |
| Dipropylene glycol  |     | 229.4 | 23.7  |
| Disiloxane, pentamethyl-•   |     |       | 4.1   |
| Dodecane  |     | 122.8 | 11.2  |

| COMPOUND IDENTIFIED                | 0.0 | 1.0   | 6.0   |
|------------------------------------|-----|-------|-------|
| Ethane, 1 , 1,1-trichloro          |     | 4.7   |       |
| Ethanol                            |     | 210.4 | 4.7   |
| Ethanol, 2-(2-butoxyethoxy)        |     | 224.8 | 105.2 |
| Ethanol, 2-butoxy                  |     | 35.6  | 3.7   |
| Ethanol, 2-phenoxy-                |     | 84.4  | 36.6  |
| Furan,tetrahydro; THF              | 2.0 |       |       |
| Heptadecane*                       |     |       | 2.4   |
| Heptanal (Heptaldehyde)*           |     |       | 2.8   |
| Heptane                            |     | 15.6  |       |
| Heptane, 2,4-dimethyl              |     | 25.0  |       |
| Heptane, 2,5-dimethyl*             |     | 13.6  |       |
| Heptane, 3-methyl                  |     | 10.9  |       |
| Heptane, 3-methylene- (9Cl)*       |     | 16.6  |       |
| Heptane, 4-methyl                  |     | 12.4  |       |
| Hexadecane (Cetane)                |     | 235.7 | 10.5  |
| Hexadecane, 2,6,10-trimethyl-*     |     |       | 10.0  |
| Hexane, 2,2,5-trimethyl- (8Cl9Cl)* |     | 126.6 | 5.2   |
| Hexane, 2,2-dimethyl               |     | 67.5  |       |
| Hexane, 3-methyl                   |     | 15.8  |       |
| Hexanoic acid                      |     | 77.6  | 66.4  |
| Isobornyl acetate*                 |     | 363.5 |       |
| Methanesulfonyl chloride*          |     |       | 1.9   |
| Morpholine, 4-methyl               | 6.6 |       | 3.4   |
| Naphthalene                        |     | 69.7  | 22.3  |
| Naphthalene, 1,2-dimethyl-*        |     | 79.8  | 3.4   |
| Naphthalene, 1,4,6-trimethyl*      |     | 120.2 | 8.2   |
| Naphthalene, 1 ,4-dimethyl*        |     | 193.8 | 5.2   |
| Naphthalene, 1,6-dimethyl*         |     | 199.9 | 10.7  |

| COMPOUND IDENTIFIED  | 0.0 | 1.0   | 6.0  |
|--|-----|-------|------|
| Naphthalene,<br><b>1,6-dimethyl-4-(1-methylethyl)-*</b>  |     | 436.3 |      |
| Naphthalene, 1-methyl  |     | 23.5  | 5.3  |
| Naphthalene, <b>2,3,6-trimethyl-</b><br><b>(8Cl9Cl)*</b>   |     |       | 18.5 |
| Naphthalene, <b>2,7-dimethyl*</b>  |     | 96.0  | 3.8  |
| Naphthalene, 2-methyl  |     | 28.3  | 10.8 |
| Naphthalene,<br><b>6(1-ethylpropyl)-1,2,3,4-tetrahydro</b><br>•  |     | 123.9 | 9.3  |
| Naphthalene,<br><b>decahydro-4a-methyl-1-methylene</b><br><b>-7-(1-methylethylidene)- (4aR-<br/>trans)-*</b> |     | 298.7 | 12.4 |
| Nonane   |     | 84.8  | 2.0  |
| Nonane, 3-methyl   |     | 73.1  |      |
| Nonane, 4-methyl   |     | 31.7  | 3.3  |
| Octane   |     | 23.1  |      |
| Octane, <b>2,6-dimethyl</b>  |     | 121.1 | 15.6 |
| Octane, 2-methyl   |     | 17.7  |      |
| Octane, 3-methyl   |     | 43.6  | 14.4 |
| Octane, <b>4-ethyl*</b>  |     | 75.8  |      |
| Pentadecane,<br><b>2,6,10,14-tetramethyl-*</b>   |     | 469.9 |      |
| Pentane, <b>2,2,4-trimethyl</b><br><b>(Isooctane)</b>  |     | 560.7 |      |
| Pentane, <b>2,3,4-trimethyl</b>  |     | 61.6  |      |
| Pentane, <b>2,3-dimethyl (8Cl9Cl)*</b>   |     | 11.8  |      |
| Pentane, <b>2,4-dimethyl*</b>  |     | 8.4   |      |
| Pentanoic acid ( <b>Valeric</b> acid)  |     | 30.1  | 20.2 |
| Phenol   |     | 183.4 | 68.7 |

| COMPOUND IDENTIFIED   | 0.0 | 1.0    | 6.0  |
|---|-----|--------|------|
| Phenol,<br>4,6-di(1,1-dimethylethyl)-2-methyl-•                     |     | 79.6   | 2.0  |
| Phenol, bis(1,1-dimethylethyl)-(9Cl)*                               |     |        | 8.0  |
| Propanedioic acid, diproyl-, diethyl ester*                         |     | 747.7  | 59.4 |
| Pyridine,<br>1,2,3,6-tetrahydro-4-[4,5-dihydroxyphenyl]-1 -methyl-* |     | 52.5   |      |
| TXIB<br>(2,2,4-Trimethyl-1,3-pentanediol diisobutyrate)             |     | 324.8  | 19.2 |
| Terbutol*   |     | 66.5   |      |
| Tetradecane   |     | 110.0  | 3.4  |
| Toluene (Methylbenzene)   |     | 289.2  |      |
| Tridecane   |     | 22.8   |      |
| Trimethylamine<br>(methanamine,N,N-dimethyl)*                       |     |        | 0.6  |
| Unidentified  |     | 1840.9 | 30.8 |
| Xylene (para and/or meta)   |     | 52.9   |      |
| Xylene, ortho   |     | 127.9  |      |
| t-2-Pentenal*   |     | 7.6    |      |
| t-5-Decene*   |     | 74.1   | 23.8 |

\*Mites best NIST/EPA/NIH library match only.

Values marked with • > symbol were outside linear range of detector. Concentration value is a lower bound only.

Individual compounds and TVOC (total volatile organic compounds) are calibrated relative to toluene.

TVOC detection limit: 0.9  $\mu\text{g}/\text{m}^3$ . Individual detection limits may vary, depending on instrument response.

Cushion **B** at Elevated Conditions (70 °C, 8% RH, high loading)

Environmental Chamber: **SC5**  
Product Loading: **1.65 m<sup>2</sup>/m<sup>3</sup>**  
Test Conditions: **1.00 ACH**  
**8.0% RH ± 2.0% RH**  
**70.0 degs C ± 2.0 degs C**  
Test Period: **06/21/95 - 06/21/95**

CHAMBER **TVOC** CONCENTRATIONS FROM 0.000 TO 6.000 HOURS

| ELAPSED EXPOSURE HOUR | Tvoc CONCENTRATION ( $\mu\text{g}/\text{m}^3$ ) |
|-----------------------|---|
| 0.000                 | 0.0   |
| 1.000                 | > 10620.9                                       |
| 6.000                 | > 10253.6                                       |

**IVOC CONCENTRATIONS ( $\mu\text{g}/\text{m}^3$ )** FROM 0.000 TO 6.000 HOURS

| COMPOUND IDENTIFIED                | 1.0    | 6.0   |
|------------------------------------|--------|-------|
| (tert-Butoxymethyl)oxirane*        | 37.2   |       |
| 1,2-Ethanediol (Ethylene glycol)   | 30.9   | 39.4  |
| 1,2-Propanediol (Propylene glycol) | 37.2   | 42.6  |
| 1,4-Dimethylpiperazine             |        | 9.5   |
| I-Acetoxy-2-propanol*              |        | 23.9  |
| I-Butanol (N-Butyl alcohol)        | 471 .0 | 134.8 |
| I-Dodecene                         | 29.3   |       |
| I-Hexanol, 2-ethyl                 | 71.4   | 80.7  |
| I-Nonene, 4,6,8-trimethyl- (9Cl)*  | 8.8    | 8.0   |
| 1-Octanol                          |        | 11.0  |
| I-Octanol, 2-butyl- (8Cl9Cl)       | 14.7   |       |
| 1-Octene, 3-ethyl-*                | 26.5   | 4.9   |

| COMPOUND IDENTIFIED  |        |        |
|--|--------|--------|
|  | 1.0    | 6.0    |
| <b>1-Pentanol, 3-methyl</b>  | 8.3    | 4.9    |
| <b>1-Phenyl-1-propyne</b>  |        | 13.5   |
| <b>1-Propanol (Propyl alcohol)</b>   | 118.9  | 19.2   |
| <b>1-Tetradecene</b>   | 7.6    |        |
| <b>1-Undecene</b>  | 35.3   |        |
| <b>1H-Indene,<br/>2,3-dihydro-1,1,3-trimethyl-*</b>                            | 7.9    |        |
| <b>1 H-Indene,<br/>2,3-dihydro-1,2-dimethyl-*</b>                              | 28.1   | 48.9   |
| <b>1 H-Indene,<br/>2,3-dihydro-1,3-dimethyl*</b>                               | 46.8   | 41.0   |
| <b>1 H-Indene,<br/>2,3-dihydro-4,7-dimethyl-*</b>                              | 29.5   | 25.0   |
| <b>1 H-Indene, 2,3-dihydro-4-methyl*</b>                                       | 69.4   | 59.5   |
| <b>1H-Phenanthro[9,10-d]imidazol-2-a<br/>mine+</b>                             | 81.1   | 119.1  |
| <b>2,2,4,4-Tetramethyloctane*</b>  |        |        |
| <b>2,2,4-Trimethyl-1,3-pentanediol<br/>monoisobutyrate (Texanol)</b>           | 58.0   | 95.9   |
| <b>2,2-Dimethyl-1-isopropyl-1,3-propan<br/>ediol monoisobutyrate (Texanol)</b> | 34.2   | 41.9   |
| <b>2,5-Cyclohexadiene-1,4-dione,<br/>2,6-bis(1,1-dimethylethyl)- (9'</b>       | 90.7 - | 134.4  |
| <b>2,6-Di-tert-butyl-4-methylphenol<br/>(BHT)</b>                              | 1823.6 | 2585.8 |
| <b>2-Ethylhexanoic acid</b>  | 664.5  | 1210.1 |
| <b>2-Pentanone, 4-hydroxy-4-methyl-<br/>(8CI9CI)*</b>                          | 447.3  | 303.9  |
| <b>2-Pentanone, 4-methyl (Methyl<br/>isobutyl ketone, MIBK)</b>                | 13.7   |        |
| <b>2-Propanol (Isopropanol)</b>  | 7.7    |        |
| <b>2-Propanol, 1,3-dichloro-</b>   | 53.9   | 41.3   |

| COMPOUND IDENTIFIED   | 1.0        | 6.0         |
|---|------------|-------------|
| 2-Propanol,<br>1-[1-methyl-2-(2-propenyoxy)ethoxy] <sup>1*</sup>                      | 29.0       |             |
| 2-Propenal*   |            | 12.3        |
| 2-Propenal, 2-methyl*   | 4.5        |             |
| 2-Pyrrolidinone, 1-methyl   |            | 14.4        |
| 2H-2,4a-Methanonaphthalene,<br>1,3,4,5,6,7-hexahydro-1,1,5,5-tetra<br>methyl-, (2S)-* | 40.1       | 37.9        |
| 3-Carene*   | 63.4       |             |
| 3-Pentanol*   | 6.8        |             |
| 3-Penten-2-one, 4-methyl- (8Cl9Cl)*   | 18.0       | 13.9        |
| 4-Isobutylmorpholine  | > 1135.3** | > 1534.1*** |
| Acetamide, N,N-dimethyl- (8Cl9Cl)   | 337.8      | 362.7       |
| Acetate, butyl  | 10.2       |             |
| Acetic acid   |            | 69.7        |
| Acetic acid, 2-ethylbutyl ester'  | 14.8       | 11.7        |
| Acetic acid, phenylmethyl ester   |            | 39.8        |
| Acetone (2-Propanone)   | 226.4      | 94.5        |
| Benzaldehyde  | 20.0       | 10.6        |
| Benzene, 1,2,4,5-tetramethyl*   |            | 37.3        |
| Benzene, 1,2,4-trimethyl  | 90.9       | 35.4        |
| Benzene, 1,4-dichloro   | 37.2       | 24.5        |
| Benzene, 1-ethyl-2,4-dimethyl-*   | 48.9       | 34.5        |
| Benzene, 1-ethyl-2-methyl<br>(2-Ethyltoluene)   | 38.9       | 9.5         |
| Benzene,<br>1-methyl-4-(1-methylethyl)<br>(p-Cymene; 4-Isopropyltoluene)              | 48.6       | 22.0        |
| Benzene, 4-ethyl-1,2-dimethyl*  |            | 21.6        |
| Benzene, ethyl  | 9.8        |             |

| COMPOUND IDENTIFIED                              | 1.0    | 6.0  |
|--|--------|------|
| Butanal (Butyraldehyde)                          | 5.2    |      |
| Butanoic acid                                    |        | 8.5  |
| Camphene   | 81.4   | 23.9 |
| Cedrene, a*                                      | 16.0   | 12.8 |
| Cyclohexane, hexyl*                              | 42.4   | 35.5 |
| Cyclohexanone                                    | 8.9    | 7.2  |
| Cyclohexene,<br>3-methyl-6-(1-methylethyldene)-* | 91.6   | 91.0 |
| Cyclohexene, 4-(1,1-dimethylethyl)*              |        | 51.4 |
| Cyclopentane,<br>1-methyl-3-(2-methylpropyl)-*   | 15.6   |      |
| Cyclopentanone*                                  | 11.7   | 16.6 |
| Cyclopentasiloxane, decamethyl*                  | 77.2   | 34.4 |
| Decane   | 114.1  | 18.0 |
| Decane, 2,2,4-trimethyl-*                        | 37.7   |      |
| Decane, 2,6,6-trimethyl-*                        | 43.2   |      |
| Decane, 2-methyl                                 | 40.8   |      |
| Decane, 3,8-dimethyl                             | 35.0   |      |
| Decane, S-methyl*                                | 73.2   |      |
| Dodecane   | 80.0 - | 57.4 |
| Dodecane, 2,6,10-trimethyl-*                     | 10.6   |      |
| Dodecane, 3-methyl- (8Cl9Cl)                     | 17.5   | 17.2 |
| Ethane, 1,1,1-trichloro                          | 3.1    |      |
| Ethane, 1,2-dichloro                             | 4.6    | 5.5  |
| Ethanol  | 36.3   | 6.3  |
| Ethanol, 2-(2-butoxyethoxy)                      |        | 46.3 |
| Ethanol, 2-butoxy                                | 33.4   | 23.2 |
| Ethanol, 2-chloro-*                              | 2.8    |      |

| COMPOUND IDENTIFIED  | 1.0   | 6.0   |
|--|-------|-------|
| Ethene, 1,1,2,2-tetrachloro<br>(Tetrachloroethylene)                                   | 6.0   |       |
| Ethyl<br>1-methylcyclopropanecarboxylate*  | 60.2  |       |
| Heptane, 2,2,6,6-tetramethyl   | 8.1   |       |
| Heptane, 2,4,6-trimethyl   | 8.9   |       |
| Hexane, 2,2,5-trimethyl- (8Cl9Cl)*   | 10.2  |       |
| Hexanoic acid  |       | 27.5  |
| Isobomyl acetate*  | 848.2 | 906.5 |
| Limonene (Dipentene;<br>1-Methyl-4-(1-methylethyl)cyclohexene)                         | 224.6 | 75.9  |
| Linalool*  |       | 71.2  |
| Morpholine, 4-ethyl-*  |       | 18.0  |
| N,N-Dimethyl-2-propenamide   | 12.3  | 14.4  |
| Naphthalene  | 115.0 | 95.3  |
| Naphthalene,<br>1,2,3,5,6,8a-hexahydro-4,7-dimethyl<br>-1-(1-methylethyl)-, (1S-cis)-* | 13.8  |       |
| Naphthalene, 1,4,6-trimethyl*  | 52.4  |       |
| Naphthalene, 1,4-dimethyl  | 10.5  | 10.2  |
| Naphthalene, 1,5-dimethyl-<br>(8Cl9Cl)*  | 6.6   |       |
| Naphthalene,<br>1,6-dimethyl-4-(1-methylethyl)-*                                       | 11.6  |       |
| Naphthalene, 1-(2-propenyl)*   | 3.4   | 3.0   |
| Naphthalene, 1-methyl  | 21.8  | 24.9  |
| Naphthalene, 2,3,6-trimethyl-<br>(8Cl9Cl)*   | 50.5  |       |
| Naphthalene, 2,6-dimethyl-*  | 9.5   |       |
| Naphthalene, 2-ethyl*  |       | 5.9   |

| COMPOUND IDENTIFIED   | 1.0   | 6.0   |
|---|-------|-------|
| Naphthalene, 2-methyl   | 63.4  | 69.7  |
| Naphthalene,6-(1,1-dimethylethyl)-1,<br>,2,3,4-tetrahydro-*         | 36.4  |       |
| Nonane  | 14.4  |       |
| Nonane, 3,7-dimethyl*   | 96.5  | 24.1  |
| Nonane, 3-methyl-5-propyl-*   | 33.0  |       |
| Nonane, 4,5-dimethyl*   | 140.6 |       |
| Nonane, 4-methyl  | 32.0  |       |
| Octane, 2,2,6-trimethyl*  | 50.9  | 11.8  |
| Octane, 3-ethyl-2,7-dimethyl- (9Cl)*                                | 118.8 |       |
| Pentanal (Valeraldehyde)  | 4.1   |       |
| Pentane, 2,2,4-trimethyl (Isooctane)                                | 2.7   |       |
| Phenol  | 165.1 | 168.3 |
| Phenol,<br>2,6-bis(1,1-dimethylethyl)-4-ethyl*                      | 34.6  |       |
| Phenol, 4-t-butyl<br>(4-(1 ,1-Dimethylethyl)phenol)*                |       | 10.4  |
| Pinene, $\alpha$<br>(2,6,6-Trimethyl-bicyclo[3.1.1]hept-2<br>-ene)  | 21.7  |       |
| Propane, 2-methyl-2-nitro*  | 7.3   |       |
| Propanoic acid  |       | 12.4  |
| Propanoic acid, 2,2-dimethyl-*                                      |       | 8.1   |
| Propanoic acid, 2-methyl- (9Cl)*                                    |       | 6.1   |
| Propanoic acid,<br>2-methyl-2-[(trimethylsilyl)oxy]-,<br>trimethyl* | 10.1  | 8.2   |
| Styrene   | 19.4  | 5.5   |
| Styrene, 2,5-dimethyl*  | 31.5  | 28.9  |
| Tetradecane   | 15.3  | 5.8   |

| COMPOUND IDENTIFIED                                     | 1.0         | 6.0   |
|---|-------------|-------|
| Thiophene,<br><u>2-butyl-5-(2-methylpropyl)- (9Cl)*</u> | 17.1        |       |
| Toluene (Methylbenzene)                                 | <u>29.5</u> | 7.8   |
| Tridecane   | <u>20.5</u> | 14.3  |
| Undecane  | 186.1       | 96.4  |
| Undecane, 2,3-dimethyl-*                                | 34.1        | 25.5  |
| Undecane, 2-methyl                                      | 27.6        |       |
| Undecane, 3,9-dimethyl- (8Cl)*                          | 42.8        | 42.0  |
| Undecane, 5-methyl- (8Cl9Cl)*                           | 30.6        |       |
| Unidentified  | 374.8       | 503.3 |
| Xylene (para and/or meta)                               | 40.8        |       |
| Xylene, ortho   | 28.9        | 9.4   |

\*Indicates best NIST/EPA/NIH library match only.

"Values marked with ">" symbol were outside linear range of detector. Concentration value is a lower bound only.

Individual compounds and TVOC (total volatile organic compounds) are calibrated relative to toluene.

TVOC detection limit: 0.9 µg/m<sup>3</sup>. Individual detection limits may vary, depending on instrument response.

## **APPENDIX D**

### **INDIVIDUAL VOC CONCENTRATIONS DURING EXPOSURES TO SYNTHESIZED MIXTURES**

TABLE D-I  
CONCENTRATION DATA FOR EXPOSURE TO  
PRIME URETHANE "A" TEST MIXTURE

| Compound                         | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------------------------------|---|--|--|---|-------------------|
|                                  | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol | 6.2   | 12.3   | 12.1   | 10.2                                    | 33.9              |
| 1,4-Dimethylpiperazine           | 11.1  | 16.2   | 17.2   | 14.8                                    | 21.9              |
| 1,1,1-Trichloroethane            | 6.8   | 7.0  | 7.2'   | 7.0                                     | 2.4               |
| Adiponitrile                     | 0.3   | 0.5  | 5.8  | 2.2                                     | 143               |
| Styrene                          | 5.6   | 6.1  | 6.1  | 5.9                                     | 4.2               |
| Octamethylcyclotetrasiloxane     | 3.0   | 3.2  | 3.3  | 3.2                                     | 3.5               |
| Ether*                           |   | 0.8  | 1.3  |   |                   |
| Ethyl acetate*                   |   | 0.5  | 0.8  |   |                   |
| Methanonaphthalene*              | 0.2   | 0.3  | 0.3  |   |                   |
| Iooctane*                        |   | 0.2  | 0.2  |   |                   |
| Methanol*                        | 0.1   | 0.2  |  |   |                   |

| Compound                      | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|-------------------------------|---|--|--|---|-------------------|
|                               | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| Hexane*                       |   | 0.1  | < 0.1  |   |                   |
| 1,1-Dichloroethene*           |   | 0.1  |  |   |                   |
| Hexamethylcyclotetrasiloxane* |   | 0.1  |  |   |                   |
| Benzene*                      |   |  | 0.1  |   |                   |

\*Quantitation relative to toluene response only.

**TABLE D-2**

**CONCENTRATION DATA FOR EXPOSURE TO  
 PRIME URETHANE "A" TEST MIXTURE WITH BHT REMOVED**

| Compound                         | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------------------------------|---|--|--|---|-------------------|
|                                  | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol | nd  | nd   | nd   | nd                                      | ---               |
| 1,4-Dimethylpiperazine           | 13.8  | 19.2   | 15.8   | 16.3                                    | 16.8              |
| 1,1,1-Trichloroethane            | 10.0  | 10.5   | 8.6  | 9.7                                     | 10.0              |
| Adiponitrile                     | nd  | 0.2  | 1.1  | 0.4                                     | 132               |
| Styrene                          | 6.2   | 6.7  | 6.7  | 6.5                                     | 4.4               |
| Octamethylcyclotetrasiloxane     | 4.4   | 4.3  | 4.6  | 4.4                                     | 3.4               |
| Methanol*                        | 4.5   | 3.1  | 1.9  |   |                   |
| Pyrazine*                        | 0.1   | 0.2  | 0.3  |   |                   |
| Ether*                           |   |  | 3.5  |   |                   |
| 1,1-Dichloroethene*              |   |  | 1.7  |   |                   |
| Isooctane*                       |   |  | 0.6  |   |                   |

| Compound       | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------------|---|--|--|---|-------------------|
|                | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| Ethyl acetate* |   |  | 0.5  |   |                   |
| Acetone*       |   |  | 0.3  |   |                   |
| Hexane*        |   |  | 0.1  |   |                   |
| Benzene*       |   |  | 0.1  |   |                   |

\*Quantitation relative to toluene response only.

TABLE D-3  
CONCENTRATION DATA FOR EXPOSURE TO  
PRIME URETHANE "A" TEST MIXTURE  
WITH BHT AND 1,4-DIMETHYLPIPERAZINE REMOVED

| Compound                         | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------------------------------|--|---|---|---|---|-------------------|
|                                  | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10-30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol | nd   | nd  | nd  | nd  | nd                                      | ---               |
| 1,4-Dimethylpiperazine           | nd   | nd  | nd  | nd  | nd                                      | ---               |
| 1,1, 1-Trichloroethane           | 2.9  | 2.7   | 2.5   | 2.5   | 2.7                                     | 8.0               |
| Adiponitrile                     | nd   | 1.8   | 5.2   | 6.4   | 3.3                                     | 89                |
| Styrene                          | 4.6  | 5.1   | 4.3   | 4.9   | 4.7                                     | 7.6               |
| Octamethylcyclotetrasiloxane     | 0.3  | 0.8   | 1.0   | 1.4   | 0.9                                     | 51                |

TABLE D-4

CONCENTRATION DATA FOR EXPOSURE TO LOW CONCENTRATION PRIME URETHANE "A" TEST MIXTURE

| Compound                         | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------------------------------|---|--|--|---|-------------------|
|                                  | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol | 2.0   | 5.1  | 6.6  | 4.6                                     | 50.7              |
| 1,4-Dimethylpiperazine           | 0.6   | 2.3  | 3.2  | 2.0                                     | 66.9              |
| 1,1,1-Trichloroethane            | 0.9   | 1.3  | 1.6  | 1.3                                     | 28.4              |
| Adiponitrile                     | nd  | nd   | 1.0  | 0.4                                     | 173               |
| Styrene                          | 1.5   | 1.8  | 1.7  | 1.7                                     | 9.7               |
| Octamethylcyclotetrasiloxane     | 2.5   | 2.7  | 2.4  | 2.6                                     | 6.4               |
| Methanonaphthalene*              | 0.7   | 0.9  | 0.5  |   |                   |

\*Quantitation relative to toluene response only

TABLE D-5  
CONCENTRATION DATA FOR EXPOSURE TO  
PRIME URETHANE "B" TEST MIXTURE

| Compound                         | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration ( $\mu\text{g}/\text{m}^3$ ) | Precision (% RSD) |
|----------------------------------|---|--|--|---|-------------------|
|                                  | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol | 3.5   | 9.9  | 12.8   | 8.7   | 54.6              |
| N,N-Dimethylacrylamide           | 2.6   | 6.8  | 9.1  | 6.2   | 53.2              |
| 1,2,3-Trichloropropane           | 5.4   | 6.0  | 7.2  | 6.2   | 15.1              |
| N,N-Dimethylformamide            | 0.6   | 1.2  | 1.3  | 1.0   | 36.6              |
| N,N-Dimethylacetamide            | 0.2   | 0.9  | 1.2  | 0.8   | 71.2              |
| Octamethylcyclotetrasiloxane     | 4.6   | 4.6  | 4.9  | 4.7   | 3.5               |
| 1,3-Dichloro-2-propanol          | 0.8   | 2.0  | 3.0  | 1.9   | 58.7              |
| Methanonaphthalene*              | 1.2   | 1.7  |  |   |                   |
| 1-Methylpentylhydroperoxide*     | < 0.1   |  |  |   |                   |
| Ethyl acetate*                   |   | 0.1  |  |   |                   |
| Acetyl chloride*                 | 0.1   | 0.1  |  |   |                   |

| Compound                | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|-------------------------|---|--|--|---|-------------------|
|                         | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| 1-Chloro-2-propanone*   |   |  | 0.1  |   |                   |
| Ether*                  |   | 0.1  |  |   |                   |
| Toluene*                |   |  |  |   |                   |
| (Chloromethyl)-oxirane* |   | 0.1  | 0.2  |   |                   |
| Isooctane*              |   | < 0.1  |  |   |                   |
| Unidentified*           |   | 0.1  |  |   |                   |

\*Quantitation relative to toluene response only.

TABLE e-6

CONCENTRATION DATA FOR EXPOSURE TO  
 PRIME URETHANE "B" TEST MIXTURE WITH BHT REMOVED

| Compound                         | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------------------------------|---|--|--|---|-------------------|
|                                  | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol | nd  | NA*  | nd   | nd                                      | ---               |
| N,N-Dimethylacrylamide           | 22.9  | NA*  | 24.3   | 16.8                                    | 4.1               |
| 1,2,3-Trichloropropane           | 12.5  | NA*  | 14.2   | 7.7                                     | 9.1               |
| N,N-Dimethylformamide            | 3.7   | NA*  | 3.3  | 3.5                                     | 7.0               |
| N,N-Dimethylacetamide            | 6.4   | NA*  | 5.8  | 4.0                                     | 6.4               |
| Octamethylcyclotetrasiloxane     | 5.8   | NA*  | 5.9  | 6.7                                     | 0.7               |
| 1,3-Dichloro-2-propanol          | 9.5   | NA*  | 13.4   | 5.8                                     | 24.2              |
| (Chloromethyl)-oxirane**         | 0.5   | NA*  | 0.1  |   |                   |

\*NA = not available, sample lost during analysis.

\*\*Quantitation relative to toluene response only.

TABLE D-7  
CONCENTRATION DATA FOR EXPOSURE TO LOW CONCENTRATION PRIME URETHANE "B" TEST MIXTURE

| Compound                         | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------------------------------|---|--|--|---|-------------------|
|                                  | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol | 2.1   | 4.3  | 4.6  | 3.7                                     | 38.3              |
| N,N-Dimethylacrylamide           | 5.6   | 6.9  | 6.0  | 6.2                                     | 10.8              |
| 1,2,3-Trichloropropane           | 4.6   | 4.6  | 4.6  | 4.6                                     | 0.7               |
| N,N-Dimethylformamide            | 0.7   | 0.8  | 0.7  | 0.7                                     | 7.7               |
| N,N-Dimethylacetamide            | 0.5   | 0.8  | 0.5  | 0.6                                     | 28.8              |
| Octamethylcyclotetrasiloxane     | 4.1   | 4.1  | 4.0  | 4.1                                     | 1.2               |
| 1,3-Dichloro-2-propanol          | 1.7   | 2.7  | 3.7  | 2.7                                     | 36.8              |
| Methanonaphthalene*              | 0.6   | 0.9  | 1.0  |   |                   |
| (Chloromethyl)-oxirane*          | 0.2   | 0.2  |  |   |                   |

\*Quantitation relative to toluene response only.

TABLE a-8  
CONCENTRATION DATA FOR EXPOSURE TO  
SPONGE RUBBER TEST MIXTURE

| Compound                | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|-------------------------|---|--|--|---|-------------------|
|                         | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| N,N-Dimethylformamide   | 0.3   | 0.2  | 0.5  | 0.3                                     | 48.1              |
| N,N-Dimethylacetamide   | 0.5   | 1.0  | 1.9  | 1.1                                     | 63.9              |
| Toluene                 | 1.0   | 0.8  | 1.1  | 1.0                                     | 14.9              |
| Undecane                | 9.2   | 8.7  | 10.0   | 9.3                                     | 6.6               |
| 1,2,4-Trimethylbenzene  | 13.8  | 12.9   | 14.9   | 13.9                                    | 7.5               |
| 2-Methylnaphthalene     | 11.7  | 11.0   | 11.7   | 11.5                                    | 3.5               |
| 1,4-Dimethylnaphthalene | 9.0   | 7.3  | 7.9  | 8.1                                     | 10.9              |
| Diphenyl ether          | 2.2   | 2.4  | 2.5  | 2.4                                     | 7.9               |
| 2-Ethyltoluene*         | 0.1   | 0.1  | 0.1  |   |                   |
| Butylbenzene*           | < 0.1   | < 0.1  | < 0.1  |   |                   |

\*Quantitation relative to toluene response only.

TABLE e-9

CONCENTRATION DATA FOR EXPOSURE TO SPONGE RUBBER  
TEST MIXTURE WITH 2-METHYLNAPHTHALENE REMOVED

| Compound                | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|-------------------------|---|--|--|---|-------------------|
|                         | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| N,N-Dimethylformamide   | 0.2   | 0.5  | 0.7  | 0.5                                     | 46.5              |
| N,N-Dimethylacetamide   | 0.5   | 2.0  | 2.4  | 1.6                                     | 59.8              |
| Toluene                 | 1.0   | 1.4  | 1.5  | 1.3                                     | 22.6              |
| Undecane                | 8.3   | 10.3   | 12.1   | 10.2                                    | 18.5              |
| 1,2,4-Trimethylbenzene  | 13.3  | 16.7   | 18.7   | 16.2                                    | 17.0              |
| 2-Methylnaphthalene     | 0.2   | 0.3  | 0.5  | 0.3                                     | 52.4              |
| 1,4-Dimethylnaphthalene | 1.2   | 2.5  | 9.8  | 4.5                                     | 103               |
| Diphenyl ether          | 0.5   | 1.6  | 2.1  | 1.4                                     | 56.8              |
| 2-Ethyltoluene*         |   | 0.2  | 0.2  |   |                   |
| 1-Methylpropylbenzene*  |   |  | < 0.1  |   |                   |
| Butylbenzene*           | < 0.1   | 0.1  | 0.1  |   |                   |

\*Quantitation relative to toluene response only.

TABLE 0-C-a

SUMMARY CONCENTRATION DATA FOR EXPOSURE TO  
SPONGE RUBBER TEST MIXTURE WITH 2-METHYLNAPHTHALENE  
AND 1,4-DIMETHYLNAPHTHALENE REMOVED

| Compound                | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|-------------------------|---|--|--|---|-------------------|
|                         | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| N,N-Dimethylformamide   | nd  | nd   | nd   | nd                                      | ---               |
| N,N-Dimethylacetamide   | nd  | nd   | 0.6  | 0.2                                     | 173               |
| Toluene                 | 0.9   | 1.2  | 1.2  | 1.1                                     | 17.6              |
| Undecane                | 7.6   | 11.7   | 13.9   | 11.1                                    | 28.7              |
| 1,2,4-Trimethylbenzene  | 11.9  | 16.1   | 17.4   | 15.1                                    | 19.0              |
| 2-Methylnaphthalene     | nd  | nd   | nd   | nd                                      | ---               |
| 1,4-Dimethylnaphthalene | nd  | nd   | nd   | nd                                      | ---               |
| Diphenyl ether          | nd  | 1.0  | 2.4  | 1.1                                     | 107               |
| 2-Ethyltoluene*         | 0.1   | 0.1  | 0.2  |   |                   |

\*Quantitation relative to toluene response only.

TABLE CC-a

CONCENTRATION DATA FOR EXPOSURE TO TOW CONCENTRATION SPONGE RUBBER TEST MIXTURE  
(A)

| Compound                | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration ( $\mu\text{g}/\text{m}^3$ ) | Precision (% RSD) |
|-------------------------|--|---|---|---|---|-------------------|
|                         | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10 - 30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| N,N-Dimethylformamide   | nd   | nd  | nd  | nd  | nd  | —                 |
| N,N-Dimethylacetamide   | nd   | nd  | nd  | nd  | nd  | —                 |
| Toluene                 | 0.4  | 0.4   | 0.4   | 0.3   | 0.4   | 8.5               |
| Undecane                | 4.7  | 5.8   | 5.9   | 5.1   | 5.4   | 10.0              |
| 1,2,4-Trimethylbenzene  | 7.4  | 8.3   | 8.4   | 7.6   | 8.0   | 6.5               |
| 2-Methylnaphthalene     | 3.3  | 5.4   | 5.8   | 5.8   | 5.1   | 23.5              |
| 1,4-Dimethylnaphthalene | 2.9  | 6.6   | 9.2   | 10.5  | 7.3   | 46.1              |
| Diphenyl ether          | 0.4  | 1.1   | 1.3   | 1.5   | 1.1   | 42.5              |
| 2-Ethyltoluene*         | < 0.1  | < 0.1   | < 0.1   | < 0.1   |   |                   |
| 1-Methylnaphthalene*    | 0.1  | 0.2   | 0.2   | 0.2   |   |                   |

\*Quantitation relative to toluene response only.

TABLE 2C-2Z  
 CONCENTRATION DATA FOR EXPOSURE TO LOWER CONCENTRATION SPONGE RUBBER TEST  
 MIXTURE (B)

| Compound                | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|-------------------------|--|---|---|---|---|-------------------|
|                         | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10 - 30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| N,N-Dimethylformamide   | nd   | 0.2   | nd  | 0.1   | 0.1                                     | 0.0               |
| N,N-Dimethylacetamide   | nd   | 0.5   | nd  | 0.5   | 0.3                                     | 116               |
| Toluene                 | 0.2  | 0.5   | 0.2   | 0.3   | 0.3                                     | 42.7              |
| Undecane                | 2.4  | 4.7   | 2.6   | 3.7   | 3.4                                     | 31.8              |
| 1,2,4-Trimethylbenzene  | 9.0  | 7.2   | 4.8   | 5.7   | 9.6                                     | 19.5              |
| 2-Methylnaphthalene     | 0.9  | 3.1   | 1.2   | 2.4   | 1.9                                     | 54.2              |
| 1,4-Dimethylnaphthalene | 0.7  | 2.0   | 1.1   | 1.7   | 1.4                                     | 43.4              |
| Diphenyl ether          | 0.2  | 0.5   | 0.4   | 0.5   | 0.4                                     | 40.5              |
| 2-Ethyltoluene*         | < 0.1  | 0.1   | < 0.1   | < 0.1   |   |                   |
| 1-Methylnaphthalene*    | < 0.1  | 0.1   | < 0.1   | 0.1   |   |                   |
| 1-Undecene*             | < 0.1  | < 0.1   |   |   |   |                   |

| Compound                          | Concentration (mg/m <sup>3</sup> )         |   |   |   | Mean Concentration ( $\mu\text{g}/\text{m}^3$ ) | Precision (% RSD) |
|-----------------------------------|--|---|---|---|---|-------------------|
|                                   | Sample 1<br>(first 10 minutes of exposure) | Sample 2<br>(10 - 30 minutes of exposure) | Sample 3<br>(30 - 50 minutes of exposure) | Sample 4<br>(last 10 minutes of exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol* |  | < 0.1                                     |   |   |   |                   |
| Isooctane*                        |  |   |   | < 0.1                                     |   |                   |
| Benzene*                          | < 0.1                                      |   |   |   |   |                   |

\*Quantitation relative to toluene response only.

TABLE D-13  
 CONCENTRATION DATA FOR EXPOSURE TO LOWEST CONCENTRATION SPONGE RUBBER TEST  
 MIXTURE (C)

| Compound                      | Concentration (mg/m <sup>3</sup> )         |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|-------------------------------|--|---|---|---|---|-------------------|
|                               | Sample 1<br>(first 10 minutes of exposure) | Sample 2<br>(10 - 30 minutes of exposure) | Sample 3<br>(30 - 50 minutes of exposure) | Sample 4<br>(last 10 minutes of exposure) |   |                   |
| N,N-Dimethylformamide         | 0.7  | 0.7                                       | 0.9                                       | 0.9                                       | 0.8                                     | 11.7              |
| N,N-Dimethylacetamide         | 0.4  | 0.6                                       | 0.5                                       | 0.7                                       | 0.5                                     | 29.6              |
| Toluene                       | 0.2  | 0.1                                       | 0.1                                       | 0.1                                       | 0.1                                     | 98.0              |
| Undecane                      | 2.2  | 2.2                                       | 2.1                                       | 0.7                                       | 1.8                                     | 39.7              |
| 1,2,4-Trimethylbenzene        | 3.5  | 3.3                                       | 3.3                                       | 2.9                                       | 3.2                                     | 7.8               |
| 2-Methylnaphthalene           | 1.1  | 1.8                                       | 1.0                                       | 0.2                                       | 1.0                                     | 64.8              |
| 1,4-Dimethylnaphthalene       | 0.9  | 2.7                                       | 0.7                                       | 0.1                                       | 1.1                                     | 101               |
| Diphenyl ether                | 0.1  | 0.3                                       | 0.3                                       | 0.1                                       | 0.2                                     | 38.2              |
| Trichlorofluoromethane*       |  | 0.1                                       |   |   |   |                   |
| Octamethylcyclotetrasiloxane* | < 0.1                                      |   |   |   |   |                   |

\*Quantitation relative to toluene response only.

**TABLE D-14**  
**CONCENTRATION DATA FOR EXPOSURE TO**  
**BONDED URETHANE TEST MIXTURE**

| Compound                             | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|--------------------------------------|---|--|--|---|-------------------|
|                                      | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol     | 3.7   | 6.3  | 7.8  | 5.9                                     | 34.3              |
| Adiponitrile                         | 0.8   | 1.8  | 2.2  | 1.6                                     | 45.1              |
| Undecene                             | 12.5  | 15.5   | 17.0   | 15.0                                    | 15.3              |
| Tridecene                            | 5.2   | 8.8  | 8.3  | 7.4                                     | 25.9              |
| Decane                               | 1.5   | 1.8  | 1.9  | 1.8                                     | 13.2              |
| N,N-Dimethylacetamide                | 9.9   | 8.5  | 9.5  | 9.3                                     | 7.8               |
| TXIB                                 | 0.5   | 0.8  | 1.7  | 1.0                                     | 61.1              |
| Methanonaphthalene*                  | 3.1   | 2.7  | 2.4  |   |                   |
| 1,3,3-Trimethylbicyclo[2.2.1]heptane |   | 0.3  |  |   |                   |
| 4-Phenylcyclohexene*                 | 0.1   | 0.1  | 0.1  |   |                   |
| 1-Decene*                            |   | < 0.1  |  |   |                   |

\*Quantitation relative to toluene response only.

TABLE D-15

CONCENTRATION DATA FOR EXPOSURE TO  
BONDED URETHANE TEST MIXTURE WITH BHT REMOVED

| Compound                         | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------------------------------|---|--|--|---|-------------------|
|                                  | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol | nd  | nd   | nd   | nd                                      | ---               |
| Adiponitrile                     | 2.9   | 3.0  | 0.8  | 2.3                                     | 54.1              |
| Undecene                         | 18-E  | 21.9   | 20.8   | 20.3                                    | 8.9               |
| Tridecene                        | 11.2  | 21.9   | Z.E.Z  | 18.8                                    | 35.1              |
| Decane                           | 1.9   | 2.3  | 2.1  | 2.1                                     | 8.7               |
| N,N-Dimethylacetamide            | 6.9   | 9.4  | 8.6  | 8.3                                     | 15.0              |
| TXIB                             | 4.7   | S.Z  | 2.6  | 4.2                                     | 33.6              |
| 2,3-Dimethyl-1,4-hexadiene*      | 0.1   | 0.1  | 0.1  |   |                   |
| 1-Dodecyne*                      | 0.1   | 0.1  | 0.1  |   |                   |
| Ethyl acetate*                   |   | 0.1  |  |   |                   |
| c-1,5-Heptadiene*                |   |  | 0.1  |   |                   |

| Compound    | Concentration (mg/m <sup>3</sup> )            |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|-------------|---|--|--|---|-------------------|
|             | Sample 1<br>(first 10 minutes<br>of exposure) | Sample 2<br>(middle 10 minutes<br>of exposure) | Sample 3<br>(last 10 minutes of<br>exposure) |   |                   |
| 1-Dodecene* |   | 0.1  |  |   |                   |

\*Quantitation relative to toluene response only.

TABLE D-16  
 CONCENTRATION DATA FOR EXPOSURE TO  
 LOW CONCENTRATION BONDED URETHANE TEST MIXTURE

| Compound                         | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------------------------------|--|---|---|---|---|-------------------|
|                                  | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10 - 30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol | 2.1  | 5.0   | 7.9   | 6.5   | 5.4                                     | 46.3              |
| Adiponitrile                     | nd   | nd  | nd  | nd  | nd                                      | ---               |
| Undecene                         | .9.1   | 9.4   | 10.6  | 11.4  | 10.1                                    | 10.7              |
| Tridecene                        | 2.7  | 5.1   | 6.2   | 7.1   | 5.3                                     | 35.8              |
| Decane                           | 0.9  | 0.9   | 1.0   | 1.2   | 1.0                                     | 12.2              |
| N,N-Dimethylacetamide            | 1.1  | 2.5   | 0.7   | 3.4   | 1.9                                     | 64.2              |
| TXIB                             | nd   | 0.1   | 0.3   | 0.3   | •2                                      | 93.1              |
| Methanonaphthalene*              | 0.2  | 0.4   | 0.5   | 0.3   |   |                   |
| Norflurane*                      | 0.1  | 0.2   |   |   |   | .                 |
| Ether*                           | 0.1  |   |   |   |   |                   |
| Isooctane*                       | 0.1  |   |   |   |   |                   |

| Compound | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------|--|---|---|---|---|-------------------|
|          | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10 - 30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| Benzene  | < 0.1  |   |   |   |   |                   |

\*Quantitation relative to toluene response only.

TABLE D-17  
 CONCENTRATION DATA FOR EXPOSURE TO SBR CARPET TEST MIXTURE

| Compound                                | Concentration (mg/m <sup>3</sup> )      |  |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|---|---|--|--|--|---|-------------------|
|   | Sample 1 (first 10 minutes of exposure) | Sample 2 (10 - 30 minutes of exposure) | Sample 3 (30 - 50 minutes of exposure) | Sample 4 (last 10 minutes of exposure) |   |                   |
| 4-PCH                                   | 5.2                                     | 8.7                                    | 8.7                                    | 9.2                                    | 8.0                                     | 23.0              |
| 4-VCH                                   | 2.3                                     | 2.3                                    | 2.5                                    | 2.3                                    | 2.4                                     | 3.1               |
| Styrene                                 | 6.2                                     | 6.3                                    | 6.3                                    | 6.2                                    | 6.3                                     | 1.0               |
| Toluene                                 | 10.6                                    | 10.8                                   | 11.1                                   | 10.4                                   | 10.7                                    | 2.6               |
| Isooctane                               | 1.9                                     | 2.2                                    | 2.2                                    | 2.0                                    | 2.1                                     | 6.8               |
| Acetic acid                             | nd                                      | nd                                     | 4.8                                    | nd                                     | 1.2                                     | 200               |
| Propylbenzene                           | 2.6                                     | 2.7                                    | 2.7                                    | 2.7                                    | 2.7                                     | 1.1               |
| Dodecanol                               | nd                                      | nd                                     | nd                                     | nd                                     | nd                                      | ---               |
| Cyclohexanol                            | 3.9                                     | 4.9                                    | 5.5                                    | 5.3                                    | 4.9                                     | 14.6              |
| Undecane                                | 6.1                                     | 7.0                                    | 7.0                                    | 7.1                                    | 6.8                                     | 6.7               |
| 3-Ethenyl-4-(1-methylethyl)cyclohexene* | 0.1                                     | 0.1                                    | 0.2                                    | 0.2                                    |   |                   |
| Octamethylcyclotetrasiloxane*           | 0.1                                     |  | < 0.1                                  | 0.3                                    |   |                   |

| Compound     | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|--------------|--|---|---|---|---|-------------------|
|              | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10 - 30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| Cyclohexene* |  | 0.1   |   | 0.1   |   |                   |

\*Quantitation relative to toluene response only.

**TABLE D-18**  
**CONCENTRATION DATA FOR EXPOSURE TO SBR CARPET TEST MIXTURE WITH DECANOL**

| Compound                             | Concentration (mg/m <sup>3</sup> )      |  |  |  | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|--------------------------------------|---|--|--|--|---|-------------------|
|                                      | Sample 1 (first 10 minutes of exposure) | Sample 2 (10 - 30 minutes of exposure) | Sample 3 (30 - 50 minutes of exposure) | Sample 4 (last 10 minutes of exposure) |   |                   |
| 4-PCH                                | 6.9                                     | 9.2                                    | 8.9                                    | NA*                                    | 8.3                                     | 14.8              |
| 4-VCH                                | 2.4                                     | 2.3                                    | 1.6                                    | NA*                                    | 2.1                                     | 21.2              |
| Styrene                              | 6.1                                     | 5.8                                    | 4.4                                    | NA*                                    | 5.5                                     | 16.7              |
| Toluene                              | 10.5                                    | 9.6                                    | 7.5                                    | NA*                                    | 9.2                                     | 16.8              |
| Isooctane                            | 2.2                                     | 2.0                                    | 1.4                                    | NA*                                    | 1.9                                     | 22.7              |
| Acetic acid                          | nd                                      | 0.5                                    | nd                                     | NA*                                    | 0.2                                     | 173               |
| Propylbenzene                        | 2.7                                     | 2.6                                    | 2.1                                    | NA*                                    | 2.8                                     | 13.2              |
| Decanol                              | nd                                      | 5.5                                    | 12.6                                   | NA*                                    | 4.2                                     | 105               |
| Cyclohexanol                         | 6.0                                     | 5.4                                    | 3.7                                    | NA*                                    | 5.1                                     | 23.1              |
| Undecane                             | 6.2                                     | 6.2                                    | 7.0                                    | NA*                                    | 6.5                                     | 7.3               |
| Substituted bicyclo[2.2.1]heptanes** | 0.2                                     | 0.2                                    | 0.2                                    |  |   |                   |

| Compound                           | Concentration (mg/m <sup>3</sup> )              |  |  |  | Mean Concentration ( $\mu\text{g}/\text{m}^3$ ) | Precision (% RSD) |
|------------------------------------|---|--|--|--|---|-------------------|
|                                    | Sample 1<br>(first 10<br>minutes of<br>amsodxa) | Sample 2<br>(10 - 30<br>minutes of<br>amsodxa) | Sample 3<br>(30 - 50<br>minutes of<br>amsodxa) | Sample 4<br>(last 10<br>minutes of<br>amsodxa) |   |                   |
| 2,6-di-tert-butyl-4-methylphenol** |   |  | 0.1  |  |   |                   |
| Undecene**                         | 0.1   |  |  |  |   |                   |

\*NA = not available, sample lost during analysis.

\*\* Quantitation relative to toluene response only.

TABLE D-19  
 CONCENTRATION DATA FOR EXPOSURE TO  
 COMPLAINT SYSTEM "A" TEST MIXTURE

| Compound                         | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------------------------------|--|---|---|---|---|-------------------|
|                                  | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10 - 30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol | 0.1  | 9.0   | 15.3  | 13.6  | 9.5                                     | 71.4              |
| Phenol                           | 0.8  | 1.3   | 2.6   | 2.7   | 1.8                                     | 41.4              |
| N,N-Dimethylacetamide            | 1.5  | 2.2   | 3.8   | 2.0   | 2.4                                     | 41.3              |
| Nonanal                          | 0.6  | 0.8   | 1.4   | 1.4   | 1.1                                     | 39.8              |
| 3-Methylcyclohexanol             | 1.9  | 2.4   | 3.5   | 3.3   | 2.8                                     | 27.3              |
| 1,2,4-Trimethylbenzene           | 4.0  | 4.3   | 5.4   | 4.7   | 4.6                                     | 13.7              |
| Undecene                         | 20.1   | 21.5  | 28.5  | 26.7  | 24.2                                    | 16.7              |
| Decane                           | 4.7  | 4.5   | 5.6   | 4.8   | 4.9                                     | 10.0              |
| o-Hydroxybiphenyl                | 14.6   | < 0.1   | nd  | nd  | 3.7                                     | 199               |
| 3-Methylhexane*                  | 0.1  | 0.1   | 0.1   | 0.1   |   |                   |
| 2-Methylhexane*                  | 0.1  | 0.1   | 0.1   | 0.1   |   |                   |

| Compound                      | Concentration (mg/m <sup>3</sup> )         |   |                                       |                                       | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|-------------------------------|--|---|---------------------------------------|---------------------------------------|---|-------------------|
|                               | Sample 1<br>(first 10 minutes of exposure) | Sample 2<br>(10 - 50 minutes of exposure) | (30 Sample 503 - minutes of exposure) | (last Sample 104 minutes of exposure) |   |                   |
| Heptane*                      | 0.1  | 0.1                                       | 0.1                                   | 0.1                                   |   |                   |
| Decamethylcyclopentasiloxane* | 0.3  |   |                                       |                                       |   |                   |
| Ether*                        |  |   | 0.2                                   |                                       |   |                   |
| Methanonaphthalene*           | 0.1  |   |                                       |                                       |   |                   |
| Isooctane*                    |  | < 0.1                                     | 0.1                                   |                                       |   |                   |

\*Quantitation relative to toluene response only.

TABLE D-20  
 CONCENTRATION DATA FOR EXPOSURE TO  
 COMPLAINT SYSTEM "A" TEST MIXTURE (REPEAT)

| Compound                         | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------------------------------|--|---|---|---|---|-------------------|
|                                  | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10 - 30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol | 6.1  | 7.7   | 9.1   | 5.9   | 7.2                                     | 20.9              |
| Phenol                           | 1.7  | 2.5   | 5.0   | 4.0   | 3.3                                     | 45.5              |
| N,N-Dimethylacetamide            | 2.0  | 3.0   | 0.5   | 5.9   | 2.9                                     | 79.8              |
| Nonanal                          | 1.1  | 1.1   | 1.9   | 2.0   | 1.5                                     | 31.7              |
| 3-Methylcyclohexanol             | 3.2  | 2.4   | 3.7   | 3.8   | 3.3                                     | 19.1              |
| 1,2,4-Trimethylbenzene           | 4.5  | 3.5   | 4.9   | 4.8   | 4.4                                     | 14.0              |
| Undecene                         | > 18.2*  | > 15.5*   | > 19.8*   | > 18.7*   | > 18.0*                                 | 10.3              |
| Decane                           | 4.8  | 3.8   | 5.2   | 5.1   | 4.7                                     | 13.1              |
| 1-Undecyne**                     |  |   | < 0.1   | < 0.1   |   |                   |
| Methanonaphthalene**             |  | < 0.1   |   |   |   |                   |
| 4-Phenylcyclohexene**            | 0.1  | < 0.1   | < 0.1   |   |   |                   |

| Compound         | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|------------------|--|---|---|---|---|-------------------|
|                  | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10 - 30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| 2-Ethyltoluene** |  |   | 1.0 <   | 1.0 <   |   |                   |
| Heptane**        | 0.1  | 0.1   | 1.0   | 0.1   |   |                   |
| 2-Methylhexane** | 0.1  | 0.1   | 1.0   | 0.1   |   |                   |
| 3-Methylhexane** | 0.1  | 0.1   | 0.1   | 0.1   |   |                   |

\*Concentration estimated, data for this compound were outside of linear range of mass spec.

\*\*Quantitation relative to toluene response only.

TABLE D-21  
 CONCENTRATION DATA FOR EXPOSURE TO COMPLAINT SYSTEM "A" TEST MIXTURE  
 WITH BHT REMOVED

| Compound                         | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------------------------------|--|---|---|---|---|-------------------|
|                                  | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10 - 30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol | < 0.1  | nd  | < 0.1   | nd  | < 0.1                                   | 117               |
| Phenol                           | 1.4  | 2.3   | 1.4   | 2.9   | 2.0                                     | 37.7              |
| N,N-Dimethylacetamide            | 0.6  | 2.7   | 2.6   | 4.9   | 2.7                                     | 65.4              |
| Nonanal                          | 1.6  | 1.1   | 1.5   | 1.9   | 1.5                                     | 22.9              |
| 3-Methylcyclohexanol             | 2.5  | 2.7   | 3.4   | 3.5   | 3.0                                     | 16.4              |
| 1,2,4-Trimethylbenzene           | 4.5  | 3.9   | 4.3   | 4.8   | 4.4                                     | 7.8               |
| Undecene                         | 22.4   | 20.2  | 21.1  | 24.6  | 22.1                                    | 8.5               |
| Decane                           | 4.7  | 4.0   | 4.4   | 4.8   | 4.5                                     | 8.2               |
| o-Hydroxybiphenyl                | nd   | nd  | nd  | 0.3   | 0.1                                     | 200               |
| 1,1'-Biphenyl*                   |  | < 0.1   |   | 0.2   |   |                   |
| 2-Ethyl-1-hexanol*               |  |   |   | < 0.1   |   |                   |

| Compound               | Concentration (mg/m <sup>3</sup> )         |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|------------------------|--|---|---|---|---|-------------------|
|                        | Sample 1<br>(first 10 minutes of exposure) | Sample 2<br>(10 - 30 minutes of exposure) | Sample 3<br>(30 - 50 minutes of exposure) | Sample 4<br>(last 10 minutes of exposure) |   |                   |
| 1-Undecyne*            | < 0.1                                      | < 0.1                                     | < 0.1                                     | < 0.1                                     |   |                   |
| 5-Undecene*            |  | 0.1                                       | 0.1                                       | 0.1                                       |   |                   |
| Ethyl acetate*         | 0.1  |   |   |   |   |                   |
| Benzene*               | < 0.1                                      |   |   |   |   |                   |
| 4-Ethyltoluene*        | < 0.1                                      | < 0.1                                     | < 0.1                                     | < 0.1                                     |   |                   |
| Dibenzofuran*          |  |   |   | < 0.1                                     |   |                   |
| 1,1,1-Trichloroethane* | 0.1  | < 0.1                                     | < 0.1                                     |   |   |                   |
| Ether*                 | 0.4  |   | 0.0                                       |   |   |                   |
| 2-Eethylhexanal*       |  |   |   | 0.1                                       |   |                   |
| Isooctane*             | 0.4  | 0.2                                       | 0.2                                       |   |   |                   |
| Undecane*              | 0.2  |   |   |   |   |                   |

\*Quantitation relative to toluene response only.

TABLE a-2Z

CONCENTRATION DATA FOR EXPOSURE TO COMPLAINT SYSTEM "A" TEST MIXTURE  
WITH BHT REMOVED (REPEAT)

| Compound                         | Concentration (mg/m <sup>3</sup> )         |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|----------------------------------|--|---|---|---|---|-------------------|
|                                  | Sample 1<br>(first 10 minutes of exposure) | Sample 2<br>(10 - 30 minutes of exposure) | Sample 3<br>(30 - 50 minutes of exposure) | Sample 4<br>(last 10 minutes of exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol | nd   | nd  | nd  | nd  | nd                                      | ---               |
| Phenol                           | 0.8  | 1.4                                       | 1.6                                       | 2.0                                       | 1.5                                     | 34.1              |
| N,N-Dimethylacetamide            | 0.2  | 1.0                                       | 1.0                                       | 1.3                                       | 0.9                                     | 51.5              |
| Nonanal                          | nd   | nd  | 0.7                                       | 0.7                                       | 0.4                                     | 116               |
| 3-Methylcyclohexanol             | 1.4  | 1.7                                       | 1.6                                       | 1.8                                       | 1.7                                     | 10.6              |
| 1,2,4-Trimethylbenzene           | 4.2  | 3.8                                       | 3.8                                       | 3.9                                       | 3.9                                     | 5.4               |
| Undecene                         | 15.5                                       | 18.5                                      | 17.8                                      | 18.5                                      | 17.6                                    | 8.2               |
| Decane                           | 4.1  | 4.2                                       | 4.0                                       | 4.3                                       | 4.2                                     | 6.2               |
| 1,1'-Biphenyl*                   |  | < 0.1                                     | < 0.1                                     | 0.1                                       |   |                   |
| 4-Phenylcyclohexene*             | < 0.1                                      |   |   |   |   |                   |
| Ethyl acetate*                   |  | 0.9                                       | 0.5                                       |   |   |                   |

| Compound                    | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|-----------------------------|--|---|---|---|---|-------------------|
|                             | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10 - 30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| Benzene*                    |  | 0.1   | 0.1   |   |   |                   |
| Carbon disulfide*           |  | 0.5   | 0.4   |   |   |                   |
| Chloroform*                 |  | 0.4   | 0.2   |   |   |                   |
| Dibenzofuran*               |  | < 0.1   | 0.1   | 0.2   |   |                   |
| 1,1,1-Trichloroethane*      |  | 0.1   | 0.1   |   |   |                   |
| Ether*                      |  | 1.5   | 1.4   |   |   |                   |
| Trichloromonofluoromethane* |  | 1.1   | 1.2   |   |   |                   |
| Isooctane*                  | •  | 0.3   | 0.3   |   |   |                   |

\*Quantitation relative to toluene response only.

TABLE D-23  
 CONCENTRATION DATA FOR EXPOSURE TO COMPLAINT SYSTEM "B" TEST MIXTURE

| Compound  | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration (µg/m <sup>3</sup> ) | Precision (% RSD) |
|---|--|---|---|---|---|-------------------|
|   | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10 - 30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol                              | 3.8  | 8.4   | 9.4   | 10.1  | 7.9                                     | 36.2              |
| Isobornyl acetate   | 2.5  | 2.2   | 3.7   | 2.9   | 2.8                                     | 23.2              |
| N,N-Dimethylacetamide   | 3.2  | 2.0   | 3.7   | 2.4   | 2.8                                     | 28.5              |
| Undecene  | 3.4  | 2.4   | 3.8   | 3.1   | 3.2                                     | 18.9              |
| 4-Morpholine ethanamine                                       | 0.1  | nd  | 0.6   | nd  | 0.2                                     | 173               |
| Camphene*   | 0.1  | 0.4   | 0.2   | 0.1   |   |                   |
| Methanonaphthalene*   |  | 0.3   | 0.2   | 0.1   |   |                   |
| 6-Isopropylidene-1-methyl-<br>bicyclo[3.1.0]hexane*           |  |   | 0.1   |   |   |                   |
| cis-2-Methylene-3-(1-<br>methylethyl)cyclohexanol<br>acetate* |  | < 0.1   | < 0.1   | 0.1   |   |                   |

\*Quantitation relative to toluene response only.

TABLE D-24

CONCENTRATION DATA FOR EXPOSURE TO COMPLAINT SYSTEM "B" TEST MIXTURE WITH BHT REMOVED

| Compound                                     | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration ( $\mu\text{g}/\text{m}^3$ ) | Precision (% RSD) |
|--|--|---|---|---|---|-------------------|
|  | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10 - 30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| 2,6-Di-tert-butyl-4-methylphenol             | nd   | < 0.1   | 0.1   | 0.1   | 0.1   | 69.6              |
| Isobornyl acetate                            | 0.1  | 1.5   | 3.7   | 4.2   | 2.4   | 81.8              |
| N,N-Dimethylacetamide                        | 1.7  | 1.9   | 3.3   | 4.0   | 2.7   | 41.0              |
| Undecene                                     | 3.1  | 3.0   | 5.3   | 5.7   | 4.3   | 33.3              |
| 4-Morpholine ethanamine                      | nd   | nd  | nd  | 7.4   | 1.9   | 200               |
| Camphene*                                    | 2.1  | 0.1   | 0.1   | 0.1   |   |                   |
| 1-Dodecene*                                  |  |   |   | 0.1   |   |                   |
| 1-Octanol*                                   |  |   | 0.4   |   |   |                   |
| 4-Isobutylmorpholine*                        |  |   | 3.9   | 0.1   |   |                   |
| 1-Methyl-4-(1-methylethylidene)-cyclohexene* |  |   | 0.1   | 0.1   |   |                   |
| Quillajisol                                  |  |   | 0.1   | 0.3   |   |                   |

| Compound  | Concentration (mg/m <sup>3</sup> )               |   |   |   | Mean Concentration ( $\mu\text{g}/\text{m}^3$ ) | Precision (% RSD) |
|---|--|---|---|---|---|-------------------|
|   | Sample 1<br>(first 10<br>minutes of<br>exposure) | Sample 2<br>(10 - 30<br>minutes of<br>exposure) | Sample 3<br>(30 - 50<br>minutes of<br>exposure) | Sample 4<br>(last 10<br>minutes of<br>exposure) |   |                   |
| Octyl-oxirane*  |  |   | < 0.1   |   |   |                   |
| Isooctane*  |  |   | 0.1   |   |   |                   |
| 1,7,7-Trimethyl-<br>tricyclo[2.2.1.0 <sup>2,6</sup> ]heptane* | 0.1  |   |   |   |   |                   |
| Undecane*   |  |   | 0.4   |   |   |                   |
| 2-Methylundecane*   |  |   | 0.1   |   |   |                   |

\*Quantitation relative to toluene response only.